EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	6	(("6166849") or ("6698657") or ("4623702")).PN.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	OFF	2006/11/14 14:00
L2	6	(("6166849") or ("5698657") or ("4623702")).PN.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	OFF	2006/11/14 14:05
L3	4	Ancamine adj3 "2205"	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:05
L4	6	Ancamine adj3 "1856"	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:05

EAST Search History

L5	2	l3 and l4	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:06
L6	4	l3	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:10
L7	2	jp-61148280-$.did.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:15
L8	83951	aliphatic near5 (amin?$1 or diamin?$1 or polyamin?$1)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:16

EAST Search History

L9	75	l8 near7 (amine adj (value or number))	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:17
L10	66	l8 near7 ((glass adj transition adj temperature) or Tg)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:17
L11	1	l9 and l10	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:17
L12	21175	adhesive near7 (anchor$3 or bolt or bore or borehole)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:18

EAST Search History

L13	2	l12 and l9	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:19
L14	2	l12 and l10	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 14:19
L15	32759	amin?$1 near3 epox?$3	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 15:53
L16	1398	ratio near5 l15	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 15:53

EAST Search History

L17	13	l12 and l16	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 15:57
L18	114001	(epox?$3 or epoxidi?ed or epoxidat?$3 or epoxylat?$3) near5 (polybutadiene or novola?$1 or (phenol$2 adj resin)) or diepox?$3 or polyepox?$3 or diglycidyl$ or polyglycidyl$	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 15:59
L19	25	l9 and l18	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 16:29
L20	21	l10 and l18	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/11/14 16:29

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     (FILE 'HOME' ENTERED AT 12:48:04 ON 14 NOV 2006)

     FILE 'REGISTRY' ENTERED AT 12:48:36 ON 14 NOV 2006
L1             3 S ANCAMINE(W)1856
L2             2 S ANCAMINE(W)2205
L3             1 S ANCAMINE(W)K54
L4             1 S 90-72-2/RN

     FILE 'CAPLUS' ENTERED AT 12:54:09 ON 14 NOV 2006
L5             1 S 360771-82-0#/RN
L6             3 S 200513-58-2#/RN
L7            16 S 57214-10-5#/RN
L8             1 S L6 AND L7

     FILE 'REGISTRY' ENTERED AT 12:59:45 ON 14 NOV 2006
L9           977 S OCTYLAMINE
L10        32674 S PROPANEDIAMINE
L11        24191 S 1(W)3(W)PROPANEDIAMINE
L12            2 S NEOPENTANEDIAMINE
L13         1192 S DIMETHYL(W)1(W)3(W)PROPANEDIAMINE
L14         2049 S HEXAMETHYLENEDIAMINE
L15         2838 S DIETHYLENETRIAMINE
L16           48 S BIS(3W)AMINOPROPYL(W)AMINE
L17           10 S N(W)N(W)BIS(3W)AMINOPROPYL(W)METHYLAMINE
L18         1319 S TRIETHYLENETETRAMINE
L19            1 S 112-24-3/RN
L20          509 S TETRAETHYLENEPENTAMINE
L21            0 S 500-509
L22          509 S TETRAETHYLENEPENTAMINE
L23          132 S PENTAETHYLENEHEXAMINE
L24           19 S TRIMETHYLHEXANE(3W)DIAMINE
L25        50313 S EP/PCT

     FILE 'CAPLUS' ENTERED AT 13:10:34 ON 14 NOV 2006
L26        32989 S (111-86-4# OR 107-45-9# OR 109-76-2# OR 7328-91-8# OR 124-09-
L27         8828 S (25085-99-8# OR 1675-54-3# OR 26068-38-6#)/RN
L28        29437 S 25068-38-6#/RN
L29        79119 S L25 OR L27 OR L28
L30         1246 S 90-72-2#/RN
L31           76 S L26 AND L29 AND L30

     FILE 'STNGUIDE' ENTERED AT 13:14:47 ON 14 NOV 2006

     FILE 'CAPLUS' ENTERED AT 13:14:52 ON 14 NOV 2006

     FILE 'STNGUIDE' ENTERED AT 13:15:15 ON 14 NOV 2006

     FILE 'CA' ENTERED AT 13:37:27 ON 14 NOV 2006
L32            0 S 1971:125287/AN
L33            0 S CA71:125287W
L34            1 S 125287W

     FILE 'CAPLUS' ENTERED AT 13:38:52 ON 14 NOV 2006
L35            1 S 1971:125287#/AN
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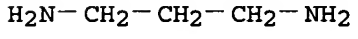
RN 111-86-4 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1-Octanamine (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Octylamine (8CI)
OTHER NAMES:
CN 1-Aminooctane
CN 1-Octylamine
CN Amine OD
CN Armeen 8
CN Armeen 8D
CN Caprylamine
CN Caprylylamine
CN Farmin 08D
CN Genamin 8R
CN Monooctylamine
CN n-Octylamine
CN NSC 9824
CN Octanamine
DR 191113-89-0
MF C8 H19 N
CI COM
LC STN Files: AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MSDS-OHS, NAPRALERT, PIRA, PS, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

$H_2N-(CH_2)_7-Me$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

4178 REFERENCES IN FILE CA (1907 TO DATE)
489 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
4185 REFERENCES IN FILE CAPLUS (1907 TO DATE)
48 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

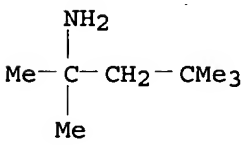
RN 109-76-2 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,3-Propanediamine (6CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN α,ω-Propanediamine
CN 1,3-Diamino-n-propane
CN 1,3-Diaminopropane
CN 1,3-Propylenediamine
CN 1,3-Trimethylenediamine
CN 3-Aminopropylamine
CN DAP
CN NSC 8154
CN TMEDA
CN Trimethylenediamine
DR 54018-94-9
MF C3 H10 N2
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MSDS-OHS, NAPRALERT, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

6193 REFERENCES IN FILE CA (1907 TO DATE)
934 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
6200 REFERENCES IN FILE CAPLUS (1907 TO DATE)
9 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

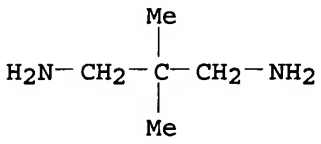
RN 107-45-9 REGISTRY
ED Entered STN: 16 Nov 1984
CN 2-Pentanamine, 2,4,4-trimethyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Butylamine, 1,1,3,3-tetramethyl- (6CI, 7CI, 8CI)
OTHER NAMES:
CN 1,1,3,3-Tetramethylbutanamine
CN 1,1,3,3-Tetramethylbutylamine
CN 2,4,4-Trimethyl-2-pentanamine
CN 2,4,4-Trimethyl-2-pentylamine
CN 2-Amino-2,4,4-trimethylpentane
CN NSC 33852
CN Primene TOA
CN tert-Octanamine
CN tert-Octylamine
DR 77658-17-4, 89961-29-5, 38724-98-0
MF C8 H19 N
CI COM
LC STN Files: ANABSTR, AQUIRE, BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DETHERM*, GMELIN*, IFICDB, IFIPAT, IFIUDB, MSDS-OHS, RTECS*, SPECINFO, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

433 REFERENCES IN FILE CA (1907 TO DATE)
31 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
433 REFERENCES IN FILE CAPLUS (1907 TO DATE)
22 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

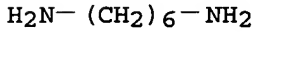
RN 7328-91-8 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,3-Propanediamine, 2,2-dimethyl- (6CI, 7CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN 1,2-Diamino-2,2-dimethylpropane
CN 1,3-Diamino-2,2-dimethylpropane
CN 2,2-Dimethyl-1,3-diaminopropane
CN 2,2-Dimethyl-1,3-propanediamine
CN 2,2-Dimethyl-1,3-propylenediamine
CN 2,2-Dimethyltrimethylenediamine
CN NSC 17719
MF C5 H14 N2
CI COM
LC STN Files: BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CSCHEM, CSNB, GMELIN*, IFICDB, IFIPAT, IFIUDB, SPECINFO, SYNTHLINE, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINECS**, NDSL**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

362 REFERENCES IN FILE CA (1907 TO DATE)
21 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
362 REFERENCES IN FILE CAPLUS (1907 TO DATE)
4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 124-09-4 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,6-Hexanediamine (7CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN α,ω-Hexanediamine
CN 1,6-Diamino-n-hexane
CN 1,6-Diaminohexane
CN 1,6-Hexylenediamine
CN 80HMD
CN Hexamethylenediamine
CN Hexylenediamine
CN Hi Perm
CN HMDA
CN NSC 9257
CN V 1
MF C6 H16 N2
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, PIRA, PROMT, PS, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

8733 REFERENCES IN FILE CA (1907 TO DATE)
1791 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
8750 REFERENCES IN FILE CAPLUS (1907 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 111-40-0 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,2-Ethanediamine, N-(2-aminoethyl)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Diethylenetriamine (8CI)
OTHER NAMES:
CN 1,4,7-Triazaheptane
CN 1,5-Diamino-3-azapentane
CN 2,2'-Diaminodiethylamine
CN 2,2'-Iminobis(ethanamine)
CN 2-(2-Aminoethylamino)ethylamine
CN 3-Azapentane-1,5-diamine
CN Ancamine DETA
CN Bis(β-aminoethyl)amine
CN Bis(2-aminoethyl)amine
CN ChS-P 1
CN DEH 20
CN DETA
CN Epicure T
CN Epon 3223
CN H 9506
CN N,N-Bis(2-aminoethyl)amine
CN N-(2-Aminoethyl)-1,2-ethanediamine
CN N-(2-Aminoethyl)ethylenediamine
CN NCI 138881
CN NSC 446
DR 859039-00-2, 8076-55-9, 53303-76-7, 54018-92-7, 59135-90-9, 94700-17-1, 98824-35-2, 73989-30-7, 26915-78-6, 419553-44-9
MF C4 H13 N3
CI COM
LC STN Files: AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MSDS-OHS, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL, VTB
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Other Sources: DSL**, EINECS**, TSCA**
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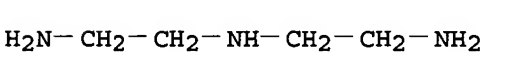
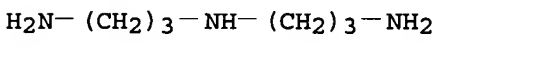


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

9317 REFERENCES IN FILE CA (1907 TO DATE)
3112 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
9332 REFERENCES IN FILE CAPLUS (1907 TO DATE)
168 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

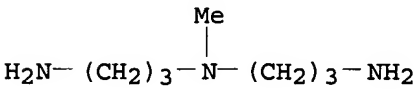
RN 56-18-8 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,3-Propanediamine, N-(3-aminopropyl)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Dipropylamine, 3,3'-diamino- (6CI, 8CI)
OTHER NAMES:
CN 1,5,9-Triazanonane
CN 1,7-Diamino-4-azaheptane
CN 1-Propanamine, 3,3'-iminobis-
CN 3,3'-Diaminodipropylamine
CN 3,3'-Iminobis(propylamine)
CN 3,3'-Iminodi(propylamine)
CN 4-Aza-1,7-diaminoheptane
CN 4-Azaheptamethylenediamine
CN 4-Azaheptane-1,7-diamine
CN Bis(3-aminopropyl)amine
CN Caldine
CN Di(3-aminopropyl)amine
CN Dipropylenetriamine
CN N-(3-Aminopropyl)-1,3-propanediamine
CN N-3-Aminopropyl-1,3-diaminopropane
CN Norspermidine
CN NSC 7773
CN P 2 (hardener)
CN sym-Norspermidine
MF C6 H17 N3
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, IFICDB, IFIPAT, IFIUDB, MEDLINE, MSDS-OHS, NAPRALERT, PHAR, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1618 REFERENCES IN FILE CA (1907 TO DATE)
352 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1618 REFERENCES IN FILE CAPLUS (1907 TO DATE)
48 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

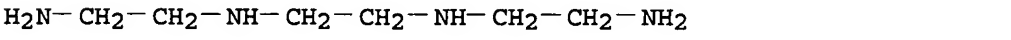
RN 105-83-9 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,3-Propanediamine, N-(3-aminopropyl)-N-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Dipropylamine, 3,3'-diamino-N-methyl- (6CI, 7CI, 8CI)
OTHER NAMES:
CN 1,7-Diamino-4-aza(methyl)heptane
CN 1,7-Diamino-4-methyl-4-azaheptane
CN 3,3'-Diamino-N-methyldipropylamine
CN 3,3'-Diamino-N-methyldipropylmine
CN 3,3'-Methyliminobispropylamine
CN 4-Aza-4-methylheptane-1,7-diamine
CN 5-Methyl-1,5,9-triazanonane
CN 5-Methyldipropylenetriamine
CN Bis(γ-aminopropyl)methylamine
CN Bis(ω-aminopropyl)methylamine
CN Bis(3-aminopropyl)methylamine
CN Di(γ-aminopropyl)methylamine
CN Methylbis(3-aminopropyl)amine
CN Methyliminobispropylamine
CN N'-Methyl-3,3'-diaminodipropylamine
CN N'-Methyldipropylenetriamine
CN N,N'-Bis(3-aminopropyl)methylamine
CN N,N-Bis(γ-aminopropyl)methylamine
CN N,N-Bis(3-aminopropyl)methylamine
CN N-(3-Aminopropyl)-N-methyl-1,3-propanediamine
CN N-Methyl-N,N-bis(3-aminopropyl)amine
CN N-Methyliminobis[1-propanamine]
CN N-Methyliminobis[propylamine]
CN NSC 8173
DR 113277-73-9, 92213-62-2
MF C7 H19 N3
CI COM
LC STN Files: BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, CSCHEM, GMELIN*, IFICDB, IFIPAT, IFIUDB, MSDS-OHS, RTECS*, SPECINFO, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

599 REFERENCES IN FILE CA (1907 TO DATE)
105 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
600 REFERENCES IN FILE CAPLUS (1907 TO DATE)
15 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

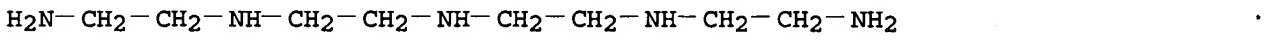
RN 112-24-3 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,2-Ethanediamine, N,N'-bis(2-aminoethyl)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Triethylenetetramine (8CI)
OTHER NAMES:
CN 1,4,7,10-Tetraazadecane
CN 1,8-Diamino-3,6-diazaoctane
CN 3,6-Diazaoctane-1,8-diamine
CN Ancamine TETA
CN Araldite Hardener HY 951
CN Araldite HY 951
CN DEH 24
CN Epicure 3234
CN HY 951
CN N,N'-Bis(2-aminoethyl)-1,2-diaminoethane
CN N,N'-Bis(2-aminoethyl)-1,2-ethanediamine
CN N,N'-Bis(2-aminoethyl)ethylenediamine
CN NSC 443
CN RT 1AX
CN Rutapox VE 2896
CN TECZA
CN TETA
CN TETA (crosslinking agent)
CN Trien
CN Trientine
CN VE 2896
CN Z1
DR 801997-18-2, 14175-14-5, 105093-20-7, 71124-11-3, 39421-77-7, 110670-33-2, 193487-08-0
MF C6 H18 N4
CI COM
LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSDRUGNEWS, IMSRESEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, PIRA, PROMT, PS, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USAN, USPAT2, USPATFULL, VETU, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

5991 REFERENCES IN FILE CA (1907 TO DATE)
1716 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
5998 REFERENCES IN FILE CAPLUS (1907 TO DATE)
114 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 112-57-2 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,2-Ethanediamine, N-(2-aminoethyl)-N'-[2-[(2-aminoethyl)amino]ethyl]-
(9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Tetraethylenepentamine (8CI)
OTHER NAMES:
CN 1,11-Diamino-3,6,9-triazaundecane
CN 1,4,7,10,13-Pentaazatridecane
CN 3,6,9-Triaza-1,11-diaminoundecane
CN 3,6,9-Triazaundecane-1,11-diamine
CN Ancamine TEPA
CN DEH 26
CN N-(2-Aminoethyl)-N'-(2-((2-aminoethyl)amino)ethyl)-1,2-ethanediamine
CN NSC 88603
CN T 11509
CN TEP
CN TEPA
CN Tetraethylpentylamine
CN Tetren
DR 778611-86-2, 115254-44-9, 675120-38-4
MF C8 H23 N5
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DETHERM*, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, MEDLINE, MSDS-OHS, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER, TULSA, USPAT2, USPATFULL, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

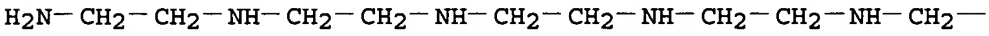


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

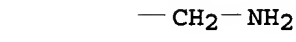
3887 REFERENCES IN FILE CA (1907 TO DATE)
1460 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
3892 REFERENCES IN FILE CAPLUS (1907 TO DATE)
96 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 4067-16-7 REGISTRY
ED Entered STN: 16 Nov 1984
CN 3,6,9,12-Tetraazatetradecane-1,14-diamine (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Pentaethylenehexamine (6CI, 7CI, 8CI)
OTHER NAMES:
CN 1,4,7,10,13,16-Hexaazahexadecane
CN 3,6,9,12-Tetraazatetradecamethylenediamine
DR 778592-37-3
MF C10 H28 N6
CI COM
LC STN Files: ANABSTR, BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, CSCHEM, IFICDB, IFIPAT, IFIUDB, MEDLINE, MSDS-OHS, RTECS*, SPECINFO, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

PAGE 1-A



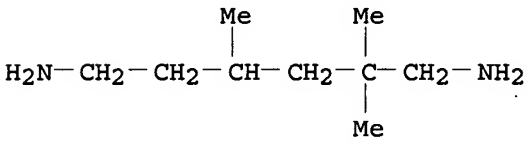
PAGE 1-B



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1070 REFERENCES IN FILE CA (1907 TO DATE)
465 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1071 REFERENCES IN FILE CAPLUS (1907 TO DATE)
13 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 3236-53-1 REGISTRY
ED Entered STN: 16 Nov 1984
CN 1,6-Hexanediamine, 2,2,4-trimethyl- (7CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN 1,6-Diamino-2,2,4-trimethylhexane
CN 2,2,4-Trimethyl-1,6-hexanediamine
CN 2,2,4-Trimethylhexamethylene-1,6-diamine
CN 2,2,4-Trimethylhexamethylenediamine
CN 2,2,4-Trimethylhexanediamine
DR 172084-55-8
MF C9 H22 N2
CI COM
LC STN Files: BEILSTEIN*, CA, CAOLD, CAPLUS, CHEMLIST, CSCHEM, IFICDB, IFIPAT, IFIUDB, MSDS-OHS, SCISEARCH, SPECINFO, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



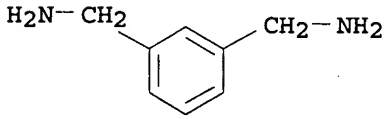
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

121 REFERENCES IN FILE CA (1907 TO DATE)
33 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
121 REFERENCES IN FILE CAPLUS (1907 TO DATE)
3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 360771-82-0 REGISTRY
ED Entered STN: 08 Oct 2001
CN Formaldehyde, polymer with 1,3-benzenedimethanamine, (chloromethyl)oxirane, 4,4'-(1-methylethylidene)bis[phenol], phenol and 2,4,6-tris[(dimethylamino)methyl]phenol (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 1,3-Benzenedimethanamine, polymer with (chloromethyl)oxirane, formaldehyde, 4,4'-(1-methylethylidene)bis[phenol], phenol and 2,4,6-tris[(dimethylamino)methyl]phenol (9CI)
CN Oxirane, (chloromethyl)-, polymer with 1,3-benzenedimethanamine, formaldehyde, 4,4'-(1-methylethylidene)bis[phenol], phenol and 2,4,6-tris[(dimethylamino)methyl]phenol (9CI)
CN Phenol, 2,4,6-tris[(dimethylamino)methyl]-, polymer with 1,3-benzenedimethanamine, (chloromethyl)oxirane, formaldehyde, 4,4'-(1-methylethylidene)bis[phenol] and phenol (9CI)
CN Phenol, 4,4'-(1-methylethylidene)bis-, polymer with 1,3-benzenedimethanamine, (chloromethyl)oxirane, formaldehyde, phenol and 2,4,6-tris[(dimethylamino)methyl]phenol (9CI)
CN Phenol, polymer with 1,3-benzenedimethanamine, (chloromethyl)oxirane, formaldehyde, 4,4'-(1-methylethylidene)bis[phenol] and 2,4,6-tris[(dimethylamino)methyl]phenol (9CI)
OTHER NAMES:
CN Ancamine 1856-Ancamine K 54-Epon 828 copolymer
MF (C15 H27 N3 O . C15 H16 O2 . C8 H12 N2 . C6 H6 O . C3 H5 Cl O . C H2 O)x
CI PMS
PCT Amino resin, Epoxy resin, Phenolic resin
SR CA
LC STN Files: CA, CAPLUS, USPATFULL

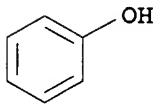
CM 1

CRN 1477-55-0
CMF C8 H12 N2



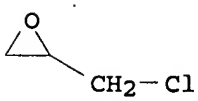
CM 2

CRN 108-95-2
CMF C6 H6 O



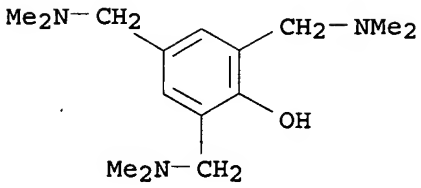
CM 3

CRN 106-89-8
CMF C3 H5 Cl O



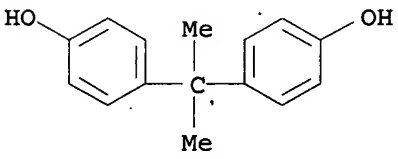
CM 4

CRN 90-72-2
CMF C15 H27 N3 O



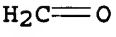
CM 5

CRN 80-05-7
CMF C15 H16 O2



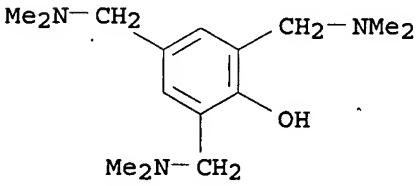
CM 6

CRN 50-00-0
CMF C H2 O



1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 90-72-2 REGISTRY
ED Entered STN: 16 Nov 1984
CN Phenol, 2,4,6-tris[(dimethylamino)methyl]- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Mesitol, α,α',α''-tris(dimethylamino)- (7CI, 8CI)
CN Mesitol, α2,α4,α6-tris(dimethylamino)- (6CI)
OTHER NAMES:
CN α,α',α''-Tris(dimethylamino)mesitol
CN 2,4,6-Tris[(dimethylamino)methyl]phenol
CN 2,4,6-Tris[(N,N-dimethylamino)methyl]phenol
CN Actiron NX 3
CN Ancamine 54
CN Ancamine K 54
CN Anchor K 54
CN Araldite DY 061
CN Araldite DY 064
CN Araldite Hardener HY 960
CN Araldite HY 960
CN Capcure EH 30
CN CK 54
CN Dabco TMR 30
CN DMF 3
CN DMP 30
CN DY 061
CN EH 30
CN Epicure 3010
CN Epilink 230
CN HY 960-1
CN K 54
CN NSC 3257
CN Protex NX 3
CN Rutapox DMP 30
CN S 41028-4
CN Sumicure D
CN TAP
CN TAP (aminophenol)
CN UP 606/2
CN Versamine CEX 13320
CN Versamine EH 30
DR 102256-77-9, 111366-64-4, 111565-20-9, 71503-41-8, 152743-49-2, 116283-46-6
MF C15 H27 N3 O
CI COM
LC STN Files: BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DETHERM*, IFICDB, IFIPAT, IFIUDB, MEDLINE, MSDS-OHS, RTECS*, SPECINFO, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1242 REFERENCES IN FILE CA (1907 TO DATE)

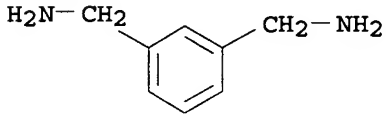
RN 200513-58-2 REGISTRY
ED Entered STN: 29 Jan 1998
CN Ancamine 2205 (9CI) (CA INDEX NAME)
ENTE An aliphatic amine (Air Products and Chemicals, Inc.)
MF Unspecified
CI COM, MAN
SR CA
LC STN Files: CA, CAPLUS, USPATFULL

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
3 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
3 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 57214-10-5 REGISTRY
ED Entered STN: 16 Nov 1984
CN Formaldehyde, polymer with 1,3-benzenedimethanamine and phenol (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 1,3-Benzenedimethanamine, polymer with formaldehyde and phenol (9CI)
CN Phenol, polymer with 1,3-benzenedimethanamine and formaldehyde (9CI)
OTHER NAMES:
CN Ancamine 1856
CN Formaldehyde-phenol-m-xylylenediamine copolymer
DR 113573-07-2
MF (C8 H12 N2 . C6 H6 O . C H2 O)x
CI PMS
PCT Amino resin, Phenolic resin
LC STN Files: CA, CAPLUS, CHEMLIST, IFICDB, IFIPAT, IFIUDB, PROMT, USPATFULL
Other Sources: TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

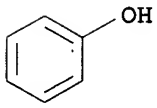
CM 1

CRN 1477-55-0
CMF C8 H12 N2



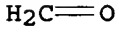
CM 2

CRN 108-95-2
CMF C6 H6 O



CM 3

CRN 50-00-0
CMF C H2 O



16 REFERENCES IN FILE CA (1907 TO DATE)
16 REFERENCES IN FILE CAPLUS (1907 TO DATE)